

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
 Data File : BM023438.D
 Acq On : 29 Oct 2019 12:21
 Operator : JU
 Sample : K5423-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE4H0

Quant Time: Oct 29 18:40:28 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 25 07:44:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	390133	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.40	136	1579818	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.27	164	956191	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.02	188	1928481	20.00	ng/ul	-0.01
78) Chrysene-d12	21.22	240	1960015	20.00	ng/ul	-0.01
86) Perylene-d12	23.41	264	2282028	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.15	96	54376	6.62	ng/uL	0.00
5) Phenol-d5	6.80	99	241996	7.39	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	6.96	67	542434	28.24	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.16	132	604246	23.59	ng/ul	-0.01
13) 4-Methylphenol-d8	8.33	113	406383	15.27	ng/ul	-0.02
19) Nitrobenzene-d5	8.77	128	353335	29.71	ng/ul	-0.02
22) 2-Nitrophenol-d4	9.49	143	383759	29.08	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.03	165	647617	24.97	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.55	131	4774	0.16	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	2293962	31.59	ng/ul	-0.01
47) Acenaphthylene-d8	13.96	160	2543615	30.58	ng/ul	-0.02
52) 4-Nitrophenol-d4	14.49	143	83026	6.60	ng/ul	0.00
58) Fluorene-d10	15.27	176	1998775	32.50	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.40	200	304924	25.40	ng/ul	-0.01
71) Anthracene-d10	17.12	188	3377668	37.06	ng/ul	-0.02
79) Pyrene-d10	19.42	212	3732347	34.63	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.27	264	4329303	37.24	ng/ul	-0.01

Target Compounds	Ovalue
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

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